

## **APPENDIX C**

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### **LABORATORY ANALYTICAL RESULTS**

[illegible]

GTEL Client ID: 7094104115110

## ANALYTICAL RESULTS

000063

Login Number: M6110437

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 06, 1996

|                |                       |             |             |             |             |
|----------------|-----------------------|-------------|-------------|-------------|-------------|
| EPA 6010A      | GTEL Sample Number    | M6110437-02 | M6110437-04 | M6110437-05 | M6110437-07 |
| Metals         | Client ID             | RP3-4       | RP3-6       | RP3-7       | RP1-4       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96    | 11/20/96    | 11/20/96    |
|                | Date Prepared         | 11/27/96    | 11/27/96    | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/04/96    | 12/04/96    | 12/04/96    | 12/04/96    |
|                | Adjustment Multiplier | 1.00        | 1.00        | 1.00        | 1.00        |
|                | Percent Solids        | 78.5        | 85.4        | 69.5        | 55.0        |

## Reporting

| Analyte  | Limit | Units | Concentration: Dry Weight |        |        |        |
|----------|-------|-------|---------------------------|--------|--------|--------|
| Barium   | 20    | mg/kg | < 20                      | < 20   | 55     | 93     |
| Cadmium  | 0.50  | mg/kg | < 0.50                    | < 0.50 | < 0.50 | < 0.51 |
| Chromium | 1.0   | mg/kg | 12                        | 3.5    | 66     | 63     |
| Lead     | 7.0   | mg/kg | < 7.0                     | < 7.0  | 25     | 31     |
| Silver   | 1.0   | mg/kg | < 1.0                     | < 1.0  | < 1.0  | < 1.0  |

|                |                       |                  |             |             |             |
|----------------|-----------------------|------------------|-------------|-------------|-------------|
| EPA 6010A      | GTEL Sample Number    | M6110437-08      | M6110437-09 | M6110437-10 | M6110437-11 |
| Metals         | Client ID             | POND 1 DUPLICATE | RP1-5       | RP2-4       | RP2-5       |
| Matrix: Solids | Date Sampled          | 11/20/96         | 11/21/96    | 11/21/96    | 11/21/96    |
|                | Date Prepared         | 11/27/96         | 11/27/96    | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/04/96         | 12/04/96    | 12/04/96    | 12/04/96    |
|                | Adjustment Multiplier | 1.05             | 1.00        | 1.00        | 1.00        |
|                | Percent Solids        | 51.9             | 62.4        | 72.9        | 80.2        |

## Reporting

| Analyte  | Limit | Units | Concentration: Dry Weight |        |        |        |
|----------|-------|-------|---------------------------|--------|--------|--------|
| Barium   | 20    | mg/kg | 95                        | 83     | 78     | 90     |
| Cadmium  | 0.50  | mg/kg | < 0.53                    | < 0.50 | < 0.50 | < 0.50 |
| Chromium | 1.0   | mg/kg | 59                        | 42     | 54     | 23     |
| Lead     | 7.0   | mg/kg | 33                        | 42     | 25     | 13     |
| Silver   | 1.0   | mg/kg | < 1.1                     | < 1.0  | < 1.0  | < 1.0  |

Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 06, 1996

#### Footnotes and Comments

##### Symbol keys (may appear beside values)

- \* - Indicates the analyte has been qualified in the narrative section of the report.
- d - Indicates the analyte was reported from a dilution other than that indicated on the report page.
- B - Organic Analyses - Indicates the analyte is found in the associated blank as well as in the sample.
- B - Inorganic Analyses - Indicates an estimated value below the EPA Contract Required Detection Limit.
- G - Indicates an estimated surrogate recovery due to dilutions.
- J - Indicates an estimated value below the reporting limit.
- U - Indicates the analyte was analyzed for but not detected.
- NA - Matrix Spike Results - Not Applicable, since the Sample Conc. exceeded four times the Spike Added.
- NA - Matrix Spike Duplicate RPD Results - Not Applicable, since the Sample Conc. exceeded four times the Spike Added.
- NA - Serial Dilution RPD Results - Not Applicable, since the Sample Conc. was less than five times the CLP Contract Required Detection Limit.

#### Inorganics

Method: EPA 6010A

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2.

Digestion is Method Specific.

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GTEL Client ID: 7094104115110

## ANALYTICAL RESULTS

Login Number: M6110437

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 06, 1996

|                |                       |             |
|----------------|-----------------------|-------------|
| EPA 6010A      | GTEL Sample Number    | M6110437-03 |
| Metals         | Client ID             | RP3-5       |
| Matrix: Solids | Date Sampled          | 11/20/96    |
|                | Date Prepared         | 11/27/96    |
|                | Date Analyzed         | 12/04/96    |
|                | Adjustment Multiplier | 1.00        |
|                | Percent Solids        | 73.5        |

| Reporting |       |       |                           |
|-----------|-------|-------|---------------------------|
| Analyte   | Limit | Units | Concentration: Dry Weight |
| Barium    | 20    | mg/kg | 27                        |
| Cadmium   | 0.50  | mg/kg | < 0.50                    |
| Chromium  | 1.0   | mg/kg | 18                        |
| Lead      | 7.0   | mg/kg | 8.9                       |
| Silver    | 1.0   | mg/kg | < 1.0                     |

Login Number: M6110437  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 05, 1996

#### Footnotes and Comments

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- J - Indicates an estimated value below the reporting limit.
- U - Indicates the analyte was analyzed for but not detected.
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- NA - Matrix Spike Duplicate RPD Results - Not Applicable, since the Sample Conc. exceeded four times the Spike Added.
- NA - Serial Dilution RPD Results - Not Applicable, since the Sample Conc. was less than five times the CLP Contract Required Detection Limit.

#### Inorganics

##### Method: EPA 7060

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2.

Digestion is Method Specific.

M6110437-03(ARSENIC): Non-Conformance Summary.

##### Matrix Spike:

MS11043703(ARSENIC): Non-Conformance Summary.

##### Method: EPA 7740

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2.

Digestion is Method Specific.

GTEL Client ID: 7094104115110

## ANALYTICAL RESULTS

Login Number: M6110437

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 05, 1996

|                |                       |             |             |             |             |
|----------------|-----------------------|-------------|-------------|-------------|-------------|
| EPA 7060       | GTEL Sample Number    | M6110437-02 | M6110437-03 | M6110437-04 | M6110437-05 |
| Metals         | Client ID             | RP3-4       | RP3-5       | RP3-6       | RP3-7       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96    | 11/20/96    | 11/20/96    |
|                | Date Prepared         | 11/27/96    | 11/27/96    | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/02/96    | 12/03/96    | 12/02/96    | 12/02/96    |
|                | Adjustment Multiplier | 1.00        | 1.50        | 1.00        | 1.00        |
|                | Percent Solids        | 78.5        | 73.5        | 85.4        | 69.5        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |     |       |
|---------|-------|-------|---------------------------|-----|-------|
| Arsenic | 1.0   | mg/kg | 3.2                       | 3.5 | < 1.0 |

|                |                       |             |                  |             |             |
|----------------|-----------------------|-------------|------------------|-------------|-------------|
| EPA 7060       | GTEL Sample Number    | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
| Metals         | Client ID             | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
|                | Date Prepared         | 11/27/96    | 11/27/96         | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/02/96    | 12/02/96         | 12/02/96    | 12/02/96    |
|                | Adjustment Multiplier | 1.00        | 1.10             | 1.00        | 1.00        |
|                | Percent Solids        | 55.0        | 51.9             | 62.4        | 72.9        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |     |     |
|---------|-------|-------|---------------------------|-----|-----|
| Arsenic | 1.0   | mg/kg | 8.3                       | 7.1 | 4.9 |

|                |                       |             |
|----------------|-----------------------|-------------|
| EPA 7060       | GTEL Sample Number    | M6110437-11 |
| Metals         | Client ID             | RP2-5       |
| Matrix: Solids | Date Sampled          | 11/21/96    |
|                | Date Prepared         | 11/27/96    |
|                | Date Analyzed         | 12/02/96    |
|                | Adjustment Multiplier | 1.00        |
|                | Percent Solids        | 80.2        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |
|---------|-------|-------|---------------------------|
| Arsenic | 1.0   | mg/kg | 2.9                       |

|                |                       |             |             |             |             |
|----------------|-----------------------|-------------|-------------|-------------|-------------|
| EPA 7740       | GTEL Sample Number    | M6110437-02 | M6110437-03 | M6110437-04 | M6110437-05 |
| Metals         | Client ID             | RP3-4       | RP3-5       | RP3-6       | RP3-7       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96    | 11/20/96    | 11/20/96    |
|                | Date Prepared         | 11/27/96    | 11/27/96    | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/03/96    | 12/03/96    | 12/03/96    | 12/03/96    |
|                | Adjustment Multiplier | 1.00        | 1.00        | 1.00        | 1.00        |
|                | Percent Solids        | 78.5        | 73.5        | 85.4        | 69.5        |

## Reporting

| Analyte  | Limit | Units | Concentration: Dry Weight |       |       |
|----------|-------|-------|---------------------------|-------|-------|
| Selenium | 1.0   | mg/kg | < 1.0                     | < 1.0 | < 1.0 |

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 05, 1996

| EPA 7740       | GTEL Sample Number    | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|----------------|-----------------------|-------------|------------------|-------------|-------------|
| Metals         | Client ID             | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
|                | Date Prepared         | 11/27/96    | 11/27/96         | 11/27/96    | 11/27/96    |
|                | Date Analyzed         | 12/03/96    | 12/03/96         | 12/03/96    | 12/03/96    |
|                | Adjustment Multiplier | 1.00        | 1.10             | 1.00        | 1.00        |
|                | Percent Solids        | 55.0        | 51.9             | 62.4        | 72.9        |

## Reporting

| Analyte  | Limit | Units | Concentration: Dry Weight |
|----------|-------|-------|---------------------------|
| Selenium | 1.0   | mg/kg | < 1.0                     |

| EPA 7740       | GTEL Sample Number    | M6110437-11 |
|----------------|-----------------------|-------------|
| Metals         | Client ID             | RP2-5       |
| Matrix: Solids | Date Sampled          | 11/21/96    |
|                | Date Prepared         | 11/27/96    |
|                | Date Analyzed         | 12/03/96    |
|                | Adjustment Multiplier | 1.00        |
|                | Percent Solids        | 80.2        |

## Reporting

| Analyte  | Limit | Units | Concentration: Dry Weight |
|----------|-------|-------|---------------------------|
| Selenium | 1.0   | mg/kg | < 1.0                     |



Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 11, 1996

#### Footnotes and Comments

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- NA - Duplicate Sample Results - Not Applicable, since the Sample Conc. was less than five times the Detection Limit.
- NA - Serial Dilution RPD Results - Not Applicable, since the Sample Conc. was less than five times the CLP Contract Required Detection Limit.

#### Inorganics

Method: EPA 7471

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds.

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 11, 1996

|                |                       |             |             |             |             |
|----------------|-----------------------|-------------|-------------|-------------|-------------|
| EPA 7471       | GTEL Sample Number    | M6110437-02 | M6110437-03 | M6110437-04 | M6110437-05 |
| Metals         | Client ID             | RP3-4       | RP3-5       | RP3-6       | RP3-7       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96    | 11/20/96    | 11/20/96    |
|                | Date Prepared         | 12/09/96    | 12/09/96    | 12/09/96    | 12/09/96    |
|                | Date Analyzed         | 12/10/96    | 12/10/96    | 12/10/96    | 12/10/96    |
|                | Adjustment Multiplier | 1.00        | 1.00        | 1.00        | 1.00        |
|                | Percent Solids        | 78.5        | 73.5        | 85.4        | 69.5        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |        |        |
|---------|-------|-------|---------------------------|--------|--------|
| Mercury | 0.25  | mg/kg | < 0.25                    | < 0.25 | < 0.25 |

|                |                       |             |                  |             |             |
|----------------|-----------------------|-------------|------------------|-------------|-------------|
| EPA 7471       | GTEL Sample Number    | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
| Metals         | Client ID             | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Matrix: Solids | Date Sampled          | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
|                | Date Prepared         | 12/09/96    | 12/09/96         | 12/09/96    | 12/09/96    |
|                | Date Analyzed         | 12/10/96    | 12/10/96         | 12/10/96    | 12/10/96    |
|                | Adjustment Multiplier | 1.00        | 1.00             | 1.00        | 1.00        |
|                | Percent Solids        | 55.0        | 51.9             | 62.4        | 72.9        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |        |        |
|---------|-------|-------|---------------------------|--------|--------|
| Mercury | 0.25  | mg/kg | < 0.25                    | < 0.25 | < 0.25 |

|                |                       |             |
|----------------|-----------------------|-------------|
| EPA 7471       | GTEL Sample Number    | M6110437-11 |
| Metals         | Client ID             | RP2-5       |
| Matrix: Solids | Date Sampled          | 11/21/96    |
|                | Date Prepared         | 12/09/96    |
|                | Date Analyzed         | 12/10/96    |
|                | Adjustment Multiplier | 1.00        |
|                | Percent Solids        | 80.2        |

## Reporting

| Analyte | Limit | Units | Concentration: Dry Weight |  |  |
|---------|-------|-------|---------------------------|--|--|
| Mercury | 0.25  | mg/kg | < 0.25                    |  |  |



**Northeast Region**

Meadowbrook Industrial Park  
Milford, NH 03055  
(603) 672-4835  
(603) 673-8105 (FAX)

January 17, 1997

Randy D. Deardorff

Buchart-Horn, Inc.  
445 W. Philadelphia St.  
York, PA 17405

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|                         |                            |
|-------------------------|----------------------------|
| RE: NEI/GTEL Client ID: | 7094104115110              |
| Login Number:           | M7010176                   |
| Project ID (number):    | 7094104115110              |
| Project ID (name):      | RFI UNIVERSITY OF MARYLAND |

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Dear Randy D. Deardorff:

Enclosed please find the analytical results for the samples received by NEI/GTEL Environmental Laboratories, Inc. on 11/25/96 under Chain-of-Custody Number(s) 64872.

A formal Quality Assurance/Quality Control (QA/QC) program is maintained by NEI/GTEL, which is designed to meet or exceed the EPA requirements. Analytical work for this project met QA/QC criteria unless otherwise stated in the footnotes. This Analytical report shall not be reproduced except in full.

GTEL is certified by the State of Maryland under certification #164.

If you have any questions regarding this analysis, or if we can be of further assistance, please call our Customer Service Representative.

Sincerely,  
NEI/GTEL Environmental Laboratories, Inc.

Susan C. Uhler  
Laboratory Director

GTEL Client ID: 7094104115110

Login Number: M7010176

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

### CONFORMANCE/NONCONFORMANCE SUMMARY

(X = Requirements Met \* = See Comments NA = Not Applicable)

| # | Conformance Item       | VOA<br>GC/MS | VOA<br>GC | SV<br>GC/MS | SV<br>GC | METALS | WET<br>CHEM |
|---|------------------------|--------------|-----------|-------------|----------|--------|-------------|
| 1 | GC/MS Tune             | —            | NA        | —           | NA       | NA     | NA          |
| 2 | Initial Calibration    | —            | —         | —           | —        | X      | —           |
| 3 | Continuing Calibration | —            | —         | —           | —        | X      | —           |
| 4 | Surrogate Recovery     | —            | —         | —           | —        | NA     | NA          |
| 5 | Holding Time           | —            | —         | —           | —        | X      | —           |
| 6 | Method Accuracy        | —            | —         | —           | —        | X      | —           |
| 7 | Method Precision       | —            | —         | —           | —        | X      | —           |
| 8 | Blank                  | —            | —         | —           | —        | X      | —           |

9 Comments: —

GTEL Client ID: 7094104115110  
Login Number: M7010176  
Project ID (Number): 7094104115110  
Project ID (Name): RFI UNIVERSITY OF MARYLAND

ANALYTICAL RESULTS

Date of Report: Jan 17, 1997

EPA 7060 GTEL Sample Number M7010176-01  
Metals Client ID FIELD BLANK  
Matrix: Aqueous Date Sampled 11/20/96  
Date Prepared 01/14/97  
Date Analyzed 01/15/97  
Adjustment Multiplier 1.00

Reporting

| Analyte | Limit | Units |
|---------|-------|-------|
| Arsenic | 10    | ug/L  |

EPA 7421 GTEL Sample Number M7010176-01  
Metals Client ID FIELD BLANK  
Matrix: Aqueous Date Sampled 11/20/96  
Date Prepared 01/14/97  
Date Analyzed 01/15/97  
Adjustment Multiplier 1.00

Reporting

| Analyte | Limit | Units |
|---------|-------|-------|
| Lead    | 4.0   | ug/L  |

EPA 7740 GTEL Sample Number M7010176-01  
Metals Client ID FIELD BLANK  
Matrix: Aqueous Date Sampled 11/20/96  
Date Prepared 01/14/97  
Date Analyzed 01/16/97  
Adjustment Multiplier 1.00

Reporting

| Analyte  | Limit | Units |
|----------|-------|-------|
| Selenium | 10    | ug/L  |

## Narrative Summary

Login Number: M7010176  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Date of Report: Jan 16, 1997

### Footnotes and Comments

#### Symbol keys (may appear beside values)

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### Inorganics

#### Method: EPA 6010A

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds.  
Digestion is Method Specific.

GTEL Client ID: 7094104115110

ANALYTICAL RESULTS

Login Number: M7010176

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Jan 16, 1997

|                 |                       |             |
|-----------------|-----------------------|-------------|
| EPA 6010A       | GTEL Sample Number    | M7010176-01 |
| Metals          | Client ID             | FIELD BLANK |
| Matrix: Aqueous | Date Sampled          | 11/20/96    |
|                 | Date Prepared         | 01/14/97    |
|                 | Date Analyzed         | 01/15/97    |
|                 | Adjustment Multiplier | 1.00        |

Reporting

| Analyte  | Limit | Units |       |
|----------|-------|-------|-------|
| Barium   | 200   | ug/L  | 200 U |
| Cadmium  | 5.0   | ug/L  | 5.0 U |
| Chromium | 10    | ug/L  | 10 U  |
| Silver   | 10    | ug/L  | 10 U  |

GTEL Milford, NH

M7010176 16:58

## PESTICIDES/PCB'S



**ANALYTICAL RESULTS**  
Organochlorine Pesticides and PCB's

GTEL Client ID: 7094104115110  
Login Number: M6110437  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
Matrix: Aqueous

|                    |             |    |    |    |
|--------------------|-------------|----|----|----|
| GTEL Sample Number | M6110437-01 | -- | -- | -- |
| Client ID          | FIELD BLANK | -- | -- | -- |
| Date Sampled       | 11/20/96    | -- | -- | -- |
| Date Prepared      | 11/25/96    | -- | -- | -- |
| Date Analyzed      | 11/26/96    | -- | -- | -- |
| Dilution Factor    | 1.00        | -- | -- | -- |

| Analyte                   | Reporting Limit | Units | Concentration: |    |    |    |
|---------------------------|-----------------|-------|----------------|----|----|----|
| PCB - Aroclor 1221        | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| PCB - Aroclor 1232        | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| PCB - Aroclor 1242 (1016) | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| PCB - Aroclor 1248        | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| PCB - Aroclor 1254        | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| PCB - Aroclor 1260        | 0.50            | ug/L  | 0.50 U         | -- | -- | -- |
| alpha-BHC                 | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| gamma-BHC(Lindane)        | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| beta-BHC                  | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Heptachlor                | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| delta-BHC                 | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Aldrin                    | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Heptachlor epoxide        | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Endosulfan I              | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| 4,4'-DDE                  | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Dieldrin                  | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| Endrin                    | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| 4,4'-DDD                  | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| Endosulfan II             | 0.05            | ug/L  | 0.05 U         | -- | -- | -- |
| 4,4'-DDT                  | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| Endrin aldehyde           | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| Endosulfan sulfate        | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| Methoxychlor              | 0.5             | ug/L  | 0.5 U          | -- | -- | -- |
| Chlordane                 | 0.1             | ug/L  | 0.1 U          | -- | -- | -- |
| Toxaphene                 | 2.0             | ug/L  | 2.0 U          | -- | -- | -- |

**Notes:**

**Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste. Physical/Chemical Methods". SW-846. Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

**M6110437-01:**

The dilution factor for pesticides = 1.00; date analyzed was 11/27/96.

GTEL Milford, NH  
M6110437

**ANALYTICAL RESULTS**  
Organochlorine Pesticides and PCB's

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
 Matrix: Solids

| GTEL Sample Number | M6110437-02 | M6110437-03 | M6110437-04 | M6110437-05 |
|--------------------|-------------|-------------|-------------|-------------|
| Client ID          | RP3-4       | RP3-5       | RP3-6       | RP3-7       |
| Date Sampled       | 11/20/96    | 11/20/96    | 11/20/96    | 11/20/96    |
| Date Prepared      | 12/04/96    | 12/04/96    | 12/04/96    | 12/04/96    |
| Date Analyzed      | 12/10/96    | 12/11/96    | 12/11/96    | 12/11/96    |
| Dilution Factor    | 1.00        | 5.00        | 1.00        | 10.0        |

| Analyte            | Reporting Limit | Units | Concentration: Dry Weight |       |       |       |
|--------------------|-----------------|-------|---------------------------|-------|-------|-------|
| alpha-BHC          | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| gamma-BHC(Lindane) | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| beta-BHC           | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Heptachlor         | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| delta-BHC          | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Aldrin             | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Heptachlor epoxide | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Endosulfan I       | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| 4,4'-DDE           | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Dieldrin           | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| Endrin             | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| 4,4'-DDD           | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| Endosulfan II      | 1.7             | ug/kg | 1.7 U                     | 8.5 U | 1.7 U | 17. U |
| 4,4'-DDT           | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| Endrin aldehyde    | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| Endosulfan sulfate | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| Methoxychlor       | 17.             | ug/kg | 17. U                     | 85. U | 17. U | 170 U |
| Chlordane          | 3.3             | ug/kg | 3.3 U                     | 16. U | 3.3 U | 33. U |
| Toxaphene          | 67.             | ug/kg | 67. U                     | 340 U | 67. U | 670 U |
| Percent Solids     | --              | %     | 78.5                      | 73.5  | 85.4  | 69.5  |

**Notes:****Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update-2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

**M6110437-03:**

Sample diluted due to non-target interference.

**M6110437-05:**

Sample diluted due to non-target interference.

**ANALYTICAL RESULTS**  
**Organochlorine Pesticides and PCB's**

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
 Matrix: Solids

| GTEL Sample Number | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|--------------------|-------------|------------------|-------------|-------------|
| Client ID          | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Date Sampled       | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
| Date Prepared      | 12/04/96    | 12/04/96         | 12/04/96    | 12/04/96    |
| Date Analyzed      | 12/11/96    | 12/11/96         | 12/11/96    | 12/11/96    |
| Dilution Factor    | 10.0        | 25.0             | 25.0        | 25.0        |

| Analyte            | Reporting Limit | Units | Concentration: Dry Weight |        |        |        |
|--------------------|-----------------|-------|---------------------------|--------|--------|--------|
| alpha-BHC          | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| gamma-BHC(Lindane) | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| beta-BHC           | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Heptachlor         | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| delta-BHC          | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Aldrin             | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Heptachlor epoxide | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Endosulfan I       | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| 4,4'-DDE           | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Dieldrin           | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| Endrin             | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| 4,4'-DDD           | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| Endosulfan II      | 1.7             | ug/kg | 17. U                     | 42. U  | 42. U  | 42. U  |
| 4,4'-DDT           | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| Endrin aldehyde    | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| Endosulfan sulfate | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| Methoxychlor       | 17.             | ug/kg | 170 U                     | 420 U  | 420 U  | 420 U  |
| Chlordane          | 3.3             | ug/kg | 33. U                     | 82. U  | 82. U  | 82. U  |
| Toxaphene          | 67.             | ug/kg | 670 U                     | 1700 U | 1700 U | 1700 U |
| Percent Solids     | --              | %     | 55.1                      | 51.9   | 62.4   | 72.9   |

**Notes:****Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update-2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

**M6110437-09:**

Sample diluted due to non-target interference.

**M6110437-10:**

Sample diluted due to non-target interference.

**ANALYTICAL RESULTS**  
**Organochlorine Pesticides and PCB's**

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
 Matrix: Solids

| GTEL Sample Number | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|--------------------|-------------|------------------|-------------|-------------|
| Client ID          | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Date Sampled       | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
| Date Prepared      | 12/04/96    | 12/04/96         | 12/04/96    | 12/04/96    |
| Date Analyzed      | 12/11/96    | 12/11/96         | 12/11/96    | 12/11/96    |
| Dilution Factor    | 1.00        | 1.00             | 1.00        | 1.00        |

| Analyte                   | Reporting Limit | Units | Concentration: Dry Weight |       |       |       |
|---------------------------|-----------------|-------|---------------------------|-------|-------|-------|
| PCB - Aroclor 1221        | 33.             | ug/kg | 33. U                     | 33. U | 33. U | 33. U |
| PCB - Aroclor 1232        | 33.             | ug/kg | 33. U                     | 33. U | 33. U | 33. U |
| PCB - Aroclor 1242 (1016) | 33.             | ug/kg | 33. U                     | 33. U | 33. U | 33. U |
| PCB - Aroclor 1248        | 33.             | ug/kg | 33. U                     | 33. U | 33. U | 33. U |
| PCB - Aroclor 1254        | 33.             | ug/kg | 33. U                     | 33. U | 33. U | 33. U |
| PCB - Aroclor 1260        | 33.             | ug/kg | 84.                       | 38.   | 68.   | 190   |
| Percent Solids            | --              | %     | 55.1                      | 51.9  | 62.4  | 72.9  |

**Notes:**

**Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

**ANALYTICAL RESULTS**  
**Organochlorine Pesticides and PCB's**

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
 Matrix: Solids

|                    |             |    |    |    |
|--------------------|-------------|----|----|----|
| GTEL Sample Number | M6110437-11 | -- | -- | -- |
| Client ID          | RP2-5       | -- | -- | -- |
| Date Sampled       | 11/21/96    | -- | -- | -- |
| Date Prepared      | 12/04/96    | -- | -- | -- |
| Date Analyzed      | 12/11/96    | -- | -- | -- |
| Dilution Factor    | 1.00        | -- | -- | -- |

| Analyte            | Reporting |       | Concentration: Dry Weight |    |    |
|--------------------|-----------|-------|---------------------------|----|----|
|                    | Limit     | Units |                           |    |    |
| alpha-BHC          | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| gamma-BHC(Lindane) | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| beta-BHC           | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Heptachlor         | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| delta-BHC          | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Aldrin             | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Heptachlor epoxide | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Endosulfan I       | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| 4,4'-DDE           | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Dieldrin           | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| Endrin             | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| 4,4'-DDD           | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| Endosulfan II      | 1.7       | ug/kg | 1.7 U                     | -- | -- |
| 4,4'-DDT           | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| Endrin aldehyde    | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| Endosulfan sulfate | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| Methoxychlor       | 17.       | ug/kg | 17. U                     | -- | -- |
| Chlordane          | 3.3       | ug/kg | 3.3 U                     | -- | -- |
| Toxaphene          | 67.       | ug/kg | 67. U                     | -- | -- |
| Percent Solids     | --        | %     | 80.2                      | -- | -- |

**Notes:****Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

**ANALYTICAL RESULTS**  
Organochlorine Pesticides and PCB's

GTEL Client ID: 7094104115110  
Login Number: M6110437  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8080  
Matrix: Solids

|                    |             |    |    |    |
|--------------------|-------------|----|----|----|
| GTEL Sample Number | M6110437-11 | -- | -- | -- |
| Client ID          | RP2-5       | -- | -- | -- |
| Date Sampled       | 11/21/96    | -- | -- | -- |
| Date Prepared      | 12/04/96    |    |    |    |
| Date Analyzed      | 12/11/96    | -- | -- | -- |
| Dilution Factor    | 1.00        | -- | -- | -- |

| Analyte                   | Reporting Limit | Units | Concentration | Dry Weight |
|---------------------------|-----------------|-------|---------------|------------|
| PCB - Aroclor 1221        | 33.             | ug/kg | 33. U         | --         |
| PCB - Aroclor 1232        | 33.             | ug/kg | 33. U         | --         |
| PCB - Aroclor 1242 (1016) | 33.             | ug/kg | 33. U         | --         |
| PCB - Aroclor 1248        | 33.             | ug/kg | 33. U         | --         |
| PCB - Aroclor 1254        | 33.             | ug/kg | 33. U         | --         |
| PCB - Aroclor 1260        | 33.             | ug/kg | 33. U         | --         |
| Percent Solids            | --              | %     | 80.2          | --         |

**Notes:**

**Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8080:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 2. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected at or above the reporting limit indicated.

# VOLATILES



000097

ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260  
 Matrix: Low Soil

| NEI/GTEL Sample Number | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|------------------------|-------------|------------------|-------------|-------------|
| Client ID              | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Date Sampled           | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
| Date Analyzed          | 12/03/96    | 12/04/96         | 12/04/96    | 12/04/96    |
| Dilution Factor        | 1.00        | 1.00             | 1.00        | 1.00        |

## Reporting

| Analyte                    | Limit | Units | Concentration: Dry Weight |       |       |       |
|----------------------------|-------|-------|---------------------------|-------|-------|-------|
| Chloromethane              | 10.   | ug/kg | 10. U                     | 10. U | 10. U | 10. U |
| Bromomethane               | 10.   | ug/kg | 10. U                     | 10. U | 10. U | 10. U |
| Vinyl chloride             | 10.   | ug/kg | 10. U                     | 10. U | 10. U | 10. U |
| Chloroethane               | 10.   | ug/kg | 10. U                     | 10. U | 10. U | 10. U |
| Methylene chloride         | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Acetone                    | 20.   | ug/kg | 59. B                     | 53. B | 62. B | 52. B |
| Carbon disulfide           | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,1-Dichloroethene         | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,1-Dichloroethane         | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,2-Dichloroethene (total) | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Chloroform                 | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,2-Dichloroethane         | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 2-Butanone                 | 20.   | ug/kg | 20. U                     | 20. U | 20. U | 20. U |
| 1,1,1-Trichloroethane      | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Carbon tetrachloride       | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Bromodichloromethane       | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,2-Dichloropropane        | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| cis-1,3-Dichloropropene    | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Trichloroethene            | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Dibromochloromethane       | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,1,2-Trichloroethane      | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Benzene                    | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| trans-1,3-Dichloropropene  | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Bromoform                  | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 4-Methyl-2-pentanone       | 20.   | ug/kg | 20. U                     | 20. U | 20. U | 20. U |
| 2-Hexanone                 | 20.   | ug/kg | 20. U                     | 20. U | 20. U | 20. U |
| Tetrachloroethene          | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| 1,1,2,2-Tetrachloroethane  | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Toluene                    | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Chlorobenzene              | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Ethylbenzene               | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Styrene                    | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Xylenes (total)            | 5.0   | ug/kg | 5.0 U                     | 5.0 U | 5.0 U | 5.0 U |
| Percent Solids             | --    | %     | 55.1                      | 51.9  | 62.4  | 72.9  |

## Notes:

## Dilution Factor:

Dilution factor indicates the adjustments made for sample dilution.

EPA 8260:  
 GTEL Milford, NH  
 M6110437

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Reissued 01/15/97D



ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110

Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260

Matrix: Low Soil

| NEI/GTEL Sample Number | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|------------------------|-------------|------------------|-------------|-------------|
| Client ID              | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Date Sampled           | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
| Date Analyzed          | 12/03/96    | 12/04/96         | 12/04/96    | 12/04/96    |
| Dilution Factor        | 1.00        | 1.00             | 1.00        | 1.00        |

| Analyte | Reporting Limit | Units | Concentration: Dry Weight |
|---------|-----------------|-------|---------------------------|
|---------|-----------------|-------|---------------------------|

Notes: (continued)

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 1. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected.

"J" indicates the presence of a compound that meets the mass spectral identification criteria, but the result is less than the reporting limit. The concentration of analytes flagged with a "J" is estimated. "B" indicates the analyte is found in the associated blank as well as the sample. It indicates possible blank contamination; The data user is warned to take appropriate action.

ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260  
 Matrix: Low Soil

|                        |             |    |    |    |
|------------------------|-------------|----|----|----|
| NEI/GTEL Sample Number | M6110437-11 | -- | -- | -- |
| Client ID              | RP2-5       | -- | -- | -- |
| Date Sampled           | 11/21/96    | -- | -- | -- |
| Date Analyzed          | 12/04/96    | -- | -- | -- |
| Dilution Factor        | 1.00        | -- | -- | -- |

| Analyte                    | Reporting Limit | Units | Concentration | Dry Weight |    |    |
|----------------------------|-----------------|-------|---------------|------------|----|----|
| Chloromethane              | 10.             | ug/kg | 10. U         | --         | -- | -- |
| Bromomethane               | 10.             | ug/kg | 10. U         | --         | -- | -- |
| Vinyl chloride             | 10.             | ug/kg | 10. U         | --         | -- | -- |
| Chloroethane               | 10.             | ug/kg | 10. U         | --         | -- | -- |
| Methylene chloride         | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Acetone                    | 20.             | ug/kg | 32. B         | --         | -- | -- |
| Carbon disulfide           | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,1-Dichloroethene         | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,1-Dichloroethane         | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,2-Dichloroethene (total) | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Chloroform                 | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,2-Dichloroethane         | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 2-Butanone                 | 20.             | ug/kg | 20. U         | --         | -- | -- |
| 1,1,1-Trichloroethane      | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Carbon tetrachloride       | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Bromodichloromethane       | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,2-Dichloropropane        | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| cis-1,3-Dichloropropene    | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Trichloroethene            | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Dibromochloromethane       | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,1,2-Trichloroethane      | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Benzene                    | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| trans-1,3-Dichloropropene  | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Bromoform                  | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 4-Methyl-2-pentanone       | 20.             | ug/kg | 20. U         | --         | -- | -- |
| 2-Hexanone                 | 20.             | ug/kg | 20. U         | --         | -- | -- |
| Tetrachloroethene          | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| 1,1,2,2-Tetrachloroethane  | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Toluene                    | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Chlorobenzene              | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Ethylbenzene               | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Styrene                    | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Xylenes (total)            | 5.0             | ug/kg | 5.0 U         | --         | -- | -- |
| Percent Solids             | --              | %     | 80.2          | --         | -- | -- |

## Notes:

## Dilution Factor:

Dilution factor indicates the adjustments made for sample dilution.

EPA 8260:

GTEL Milford, NH

M6110437

Reissued Report

ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110

Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260

Matrix: Low Soil

|                        |             |    |    |    |
|------------------------|-------------|----|----|----|
| NEI/GTEL Sample Number | M6110437-11 | -- | -- | -- |
| Client ID              | RP2-5       | -- | -- | -- |
| Date Sampled           | 11/21/96    | -- | -- | -- |
| Date Analyzed          | 12/04/96    | -- | -- | -- |
| Dilution Factor        | 1.00        | -- | -- | -- |

| Analyte | Reporting<br>Limit | Units | Concentration:Dry Weight |
|---------|--------------------|-------|--------------------------|
|---------|--------------------|-------|--------------------------|

Notes: (continued)

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 1. Analyte list may be modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected.

"J" indicates the presence of a compound that meets the mass spectral identification criteria, but the result is less than the reporting limit. The concentration of analytes flagged with a "J" is estimated. "B" indicates the analyte is found in the associated blank as well as the sample. It indicates possible blank contamination. The data user is warned to take appropriate action.

ANALYTICAL RESULTS  
Volatile Organics

GTEL Client ID: 7094104115110  
Login Number: M6110437  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260  
Matrix: Aqueous

|                    |             |             |    |    |
|--------------------|-------------|-------------|----|----|
| GTEL Sample Number | M6110437-01 | M6110437-06 | -- | -- |
| Client ID          | FIELD BLANK | TRIP BLANK  | -- | -- |
| Date Sampled       | 11/20/96    | 11/20/96    | -- | -- |
| Date Analyzed      | 12/03/96    | 12/03/96    | -- | -- |
| Dilution Factor    | 1.00        | 1.00        | -- | -- |

| Analyte                   | Reporting Limit | Units | Concentration: |       |    |
|---------------------------|-----------------|-------|----------------|-------|----|
| Dichlorodifluoromethane   | 10.             | ug/L  | 10. U          | 10. U | -- |
| Chloromethane             | 10.             | ug/L  | 10. U          | 10. U | -- |
| Vinyl chloride            | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Bromomethane              | 10.             | ug/L  | 10. U          | 10. U | -- |
| Chloroethane              | 10.             | ug/L  | 10. U          | 10. U | -- |
| Trichlorofluoromethane    | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1-Dichloroethene        | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Methylene chloride        | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| trans-1,2-Dichloroethene  | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1-Dichloroethane        | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 2,2-Dichloropropane       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| cis-1,2-Dichloroethene    | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Chloroform                | 5.0             | ug/L  | 4.1 U          | 5.0 U | -- |
| Bromochloromethane        | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1,1-Trichloroethane     | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1-Dichloropropene       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Carbon tetrachloride      | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Benzene                   | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,2-Dichloroethane        | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Trichloroethene           | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,2-Dichloropropane       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Bromodichloromethane      | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Dibromomethane            | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| cis-1,3-Dichloropropene   | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Toluene                   | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| trans-1,3-Dichloropropene | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1,2-Trichloroethane     | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,2-Dibromoethane         | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Tetrachloroethene         | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,3-Dichloropropane       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Dibromochloromethane      | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Chlorobenzene             | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Ethylbenzene              | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,1,1,2-Tetrachloroethane | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Xylenes (total)           | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,3-Dichlorobenzene       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Styrene                   | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| 1,4-Dichlorobenzene       | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |
| Bromoform                 | 5.0             | ug/L  | 5.0 U          | 5.0 U | -- |

GTEL Milford, NH  
M6110437

**ANALYTICAL RESULTS**  
**Volatile Organics**

1000102

GTEL Client ID: 7094104115110  
Login Number: M6110437  
Project ID (number): 7094104115110  
Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260  
Matrix: Aqueous

| GTEL Sample Number | M6110437-01 | M6110437-06 | -- | -- |
|--------------------|-------------|-------------|----|----|
| Client ID          | FIELD BLANK | TRIP BLANK  | -- | -- |
| Date Sampled       | 11/20/96    | 11/20/96    | -- | -- |
| Date Analyzed      | 12/03/96    | 12/03/96    | -- | -- |
| Dilution Factor    | 1.00        | 1.00        | -- | -- |

**Reporting**

| Analyte                     | Limit | Units | Concentration: |       |    |    |
|-----------------------------|-------|-------|----------------|-------|----|----|
| 1,2-Dichlorobenzene         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Isopropylbenzene            | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,1,2,2-Tetrachloroethane   | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Bromobenzene                | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2,3-Trichloropropane      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| n-Propylbenzene             | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 2-Chlorotoluene             | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,3,5-Trimethylbenzene      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 4-Chlorotoluene             | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| tert-Butylbenzene           | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2,4-Trimethylbenzene      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| sec-Butylbenzene            | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| p-Isopropyltoluene          | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| n-Butylbenzene              | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2-Dibromo-3-chloropropane | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2,4-Trichlorobenzene      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Hexachlorobutadiene         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Naphthalene                 | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2,3-Trichlorobenzene      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |

**Notes:**

**Dilution Factor:**

Dilution factor indicates the adjustments made for sample dilution.

**EPA 8260:**

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 1. "U" indicates that the analyte was analyzed for but not detected. "J" indicates the presence of a compound that meets the mass spectral identification criteria, but the result is less than the reporting limit. The concentration of analytes flagged with a "J" is estimated. "B" indicates the analyte is found in the associated blank as well as the sample. It indicates possible blank contamination: The data user is warned to take appropriate action.

ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (number): 7094104115110  
 Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260  
 Matrix: Aqueous

| NEI/GTEL Sample Number | M6110437-01 | M6110437-06 | -- | -- |
|------------------------|-------------|-------------|----|----|
| Client ID              | FIELD BLANK | TRIP BLANK  | -- | -- |
| Date Sampled           | 11/20/96    | 11/20/96    | -- | -- |
| Date Analyzed          | 12/03/96    | 12/03/96    | -- | -- |
| Dilution Factor        | 1.00        | 1.00        | -- | -- |

## Reporting

| Analyte                    | Limit | Units | Concentration: |       |    |    |
|----------------------------|-------|-------|----------------|-------|----|----|
| Chloromethane              | 10.   | ug/L  | 10. U          | 10. U | -- | -- |
| Bromomethane               | 10.   | ug/L  | 10. U          | 10. U | -- | -- |
| Vinyl chloride             | 10.   | ug/L  | 10. U          | 10. U | -- | -- |
| Chloroethane               | 10.   | ug/L  | 10. U          | 10. U | -- | -- |
| Methylene chloride         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Acetone                    | 20.   | ug/L  | 10. J          | 20. U | -- | -- |
| Carbon disulfide           | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,1-Dichloroethene         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,1-Dichloroethane         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2-Dichloroethene (total) | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Chloroform                 | 5.0   | ug/L  | 4.1 J          | 5.0 U | -- | -- |
| 1,2-Dichloroethane         | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 2-Butanone                 | 20.   | ug/L  | 20. U          | 20. U | -- | -- |
| 1,1,1-Trichloroethane      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Carbon tetrachloride       | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Bromodichloromethane       | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,2-Dichloropropane        | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| cis-1,3-Dichloropropene    | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Trichloroethene            | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Dibromochloromethane       | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,1,2-Trichloroethane      | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Benzene                    | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| trans-1,3-Dichloropropene  | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Bromoform                  | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 4-Methyl-2-pentanone       | 20.   | ug/L  | 20. U          | 20. U | -- | -- |
| 2-Hexanone                 | 20.   | ug/L  | 20. U          | 20. U | -- | -- |
| Tetrachloroethene          | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| 1,1,2,2-Tetrachloroethane  | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Toluene                    | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Chlorobenzene              | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Ethylbenzene               | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Styrene                    | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |
| Xylenes (total)            | 5.0   | ug/L  | 5.0 U          | 5.0 U | -- | -- |

## Notes:

## Dilution Factor:

Dilution factor indicates the adjustments made for sample dilution.

## EPA 8260:

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update 1. Analyte list may be  
 GTEL Milford, NH

M6110437

Page: 1

Reissued Report

ANALYTICAL RESULTS  
Volatile Organics

NEI/GTEL Client ID: 7094104115110

Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Method: EPA 8260

Matrix: Aqueous

|                        |             |             |    |    |
|------------------------|-------------|-------------|----|----|
| NEI/GTEL Sample Number | M6110437-01 | M6110437-06 | -- | -- |
| Client ID              | FIELD BLANK | TRIP BLANK  | -- | -- |
| Date Sampled           | 11/20/96    | 11/20/96    | -- | -- |
| Date Analyzed          | 12/03/96    | 12/03/96    | -- | -- |
| Dilution Factor        | 1.00        | 1.00        | -- | -- |

| Analyte | Reporting<br>Limit | Units | Concentration: |
|---------|--------------------|-------|----------------|
|---------|--------------------|-------|----------------|

Notes: (continued)

modified to include additional compounds. "U" indicates that the analyte was analyzed for but not detected.

"J" indicates the presence of a compound that meets the mass spectral identification criteria, but the result is less than the reporting limit. The concentration of analytes flagged with a "J" is estimated. "B" indicates the analyte is found in the associated blank as well as the sample. It indicates possible blank contamination; The data user is warned to take appropriate action.

# HERBICIDES



**8150 - FORM 1**  
**NYTEST ENVIRONMENTAL INC.**

**CHLORINATED PHENOXY HERBICIDES ANALYSIS DATA SHEET**

SAMPLE MATRIX : SOIL  
CONC. LEVEL : LOW  
EXTRACTION DATE : 11/29/96  
ANALYSIS DATE : 12/05/96

SAMPLE ID : RP3-4  
LAB SAMPLE ID : 2992701  
DIL FACTOR : 1.00  
% MOISTURE : 26  
INITIAL VOL. (g) : 50.01

| CMPD # | CAS Number | HERBICIDE COMPOUND | UG/KG<br>(DRY BASIS) |
|--------|------------|--------------------|----------------------|
|--------|------------|--------------------|----------------------|

|    |           |                       |       |   |
|----|-----------|-----------------------|-------|---|
| 1  | 94-75-7   | 2,4-D                 | 68    | U |
| 2  | 93-72-1   | 2,4,5-TP (Silvex)     | 6.8   | U |
| 3  | 93-76-5   | 2,4,5-T               | 6.8   | U |
| 4  | 94-82-6   | 2,4-DB                | 70    | U |
| 5  | 75-99-0   | Dalapon               | 170   | U |
| 6  | 1918-00-9 | Dicamba               | 6.8   | U |
| 7  | 88-85-7   | Dinoseb               | 32    | U |
| 8  | 120-36-5  | 2,4-DP (Dichloroprop) | 68    | U |
| 9  | 94-74-6   | MCPA                  | 27000 | U |
| 10 | 93-65-2   | MCPP                  | 27000 | U |
|    |           |                       |       |   |

000020

**8150 - FORM 1**  
**NYTEST ENVIRONMENTAL INC.**

**CHLORINATED PHENOXY HERBICIDES ANALYSIS DATA SHEET**

SAMPLE MATRIX : SOIL  
CONC. LEVEL : LOW  
EXTRACTION DATE : 11/29/96  
ANALYSIS DATE : 12/05/96

SAMPLE ID : RP3-5  
LAB SAMPLE ID : 2992702  
DIL FACTOR : 1.00  
% MOISTURE : 28  
INITIAL VOL. (g) : 50

| CMPD # | CAS Number | HERBICIDE COMPOUND | UG/KG<br>(DRY BASIS) |
|--------|------------|--------------------|----------------------|
|--------|------------|--------------------|----------------------|

|    |           |                       |       |   |
|----|-----------|-----------------------|-------|---|
| 1  | 94-75-7   | 2,4-D                 | 69    | U |
| 2  | 93-72-1   | 2,4,5-TP (Silvex)     | 6.9   | U |
| 3  | 93-76-5   | 2,4,5-T               | 6.9   | U |
| 4  | 94-82-6   | 2,4-DB                | 70    | U |
| 5  | 75-99-0   | Dalapon               | 170   | U |
| 6  | 1918-00-9 | Dicamba               | 6.9   | U |
| 7  | 88-85-7   | Dinoseb               | 33    | U |
| 8  | 120-36-5  | 2,4-DP (Dichloroprop) | 69    | U |
| 9  | 94-74-6   | MCPA                  | 28000 | U |
| 10 | 93-65-2   | MCPP                  | 28000 | U |
|    |           |                       |       |   |

000027

**8150 - FORM 1**  
**NYTEST ENVIRONMENTAL INC.**

**CHLORINATED PHENOXY HERBICIDES ANALYSIS DATA SHEET**

SAMPLE MATRIX : SOIL  
CONC. LEVEL : LOW  
EXTRACTION DATE : 11/29/96  
ANALYSIS DATE : 12/05/96

SAMPLE ID : RP-3-6  
LAB SAMPLE ID : 2992703  
DIL FACTOR : 1.00  
% MOISTURE : 20  
INITIAL VOL. (g) : 50.02

| CMPD # | CAS Number | HERBICIDE COMPOUND | UG/KG<br>(DRY BASIS) |
|--------|------------|--------------------|----------------------|
|--------|------------|--------------------|----------------------|

|    |           |                       |       |   |
|----|-----------|-----------------------|-------|---|
| 1  | 94-75-7   | 2,4-D                 | 62    | U |
| 2  | 93-72-1   | 2,4,5-TP (Silvex)     | 6.2   | U |
| 3  | 93-76-5   | 2,4,5-T               | 6.2   | U |
| 4  | 94-82-6   | 2,4-DB                | 60    | U |
| 5  | 75-99-0   | Dalapon               | 150   | U |
| 6  | 1918-00-9 | Dicamba               | 6.2   | U |
| 7  | 88-85-7   | Dinoseb               | 30    | U |
| 8  | 120-36-5  | 2,4-DP (Dichloroprop) | 62    | U |
| 9  | 94-74-6   | MCPA                  | 25000 | U |
| 10 | 93-65-2   | MCPD                  | 25000 | U |
|    |           |                       |       |   |

**000035**

**8150 - FORM 1**  
**NYTEST ENVIRONMENTAL INC.**

**CHLORINATED PHENOXY HERBICIDES ANALYSIS DATA SHEET**

SAMPLE MATRIX : SOIL  
CONC. LEVEL : LOW  
EXTRACTION DATE : 11/29/96  
ANALYSIS DATE : 12/05/96

SAMPLE ID : RP3-7  
LAB SAMPLE ID 2992704  
DIL FACTOR : 1.00  
% MOISTURE : 30  
INITIAL VOL. (g) : 50.02

| CMPD # | CAS Number | HERBICIDE COMPOUND | UG/KG<br>(DRY BASIS) |
|--------|------------|--------------------|----------------------|
|--------|------------|--------------------|----------------------|

|    |           |                       |       |   |
|----|-----------|-----------------------|-------|---|
| 1  | 94-75-7   | 2,4-D                 | 71    | U |
| 2  | 93-72-1   | 2,4,5-TP (Silvex)     | 7.1   | U |
| 3  | 93-76-5   | 2,4,5-T               | 7.1   | U |
| 4  | 94-82-6   | 2,4-DB                | 70    | U |
| 5  | 75-99-0   | Dalapon               | 180   | U |
| 6  | 1918-00-9 | Dicamba               | 7.1   | U |
| 7  | 88-85-7   | Dinoseb               | 34    | U |
| 8  | 120-36-5  | 2,4-DP (Dichloroprop) | 71    | U |
| 9  | 94-74-6   | MCPA                  | 29000 | U |
| 10 | 93-65-2   | MCP                   | 29000 | U |
|    |           |                       |       |   |

**000043**

## SEMIVOLATILES

GTEL Client ID: 7094104115110

## ANALYTICAL RESULTS

Login Number: M6110437

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 18, 1996

|                       |                       |             |
|-----------------------|-----------------------|-------------|
| EPA 8270B             | GTEL Sample Number    | M6110437-01 |
| Semivolatile Organics | Client ID             | FIELD BLANK |
| Matrix: Aqueous       | Date Sampled          | 11/20/96    |
|                       | Date Prepared         | 11/26/96    |
|                       | Date Analyzed         | 12/03/96    |
|                       | Adjustment Multiplier | 1.00        |

## Reporting

| Analyte                      | Limit | Units |      |
|------------------------------|-------|-------|------|
| Phenol                       | 10    | ug/L  | 10 U |
| bis(2-Chloroethyl) ether     | 10    | ug/L  | 10 U |
| 2-Chlorophenol               | 10    | ug/L  | 10 U |
| 1,3-Dichlorobenzene          | 10    | ug/L  | 10 U |
| 1,4-Dichlorobenzene          | 10    | ug/L  | 10 U |
| Benzyl alcohol               | 20    | ug/L  | 20 U |
| 1,2-Dichlorobenzene          | 10    | ug/L  | 10 U |
| 2-Methylphenol               | 10    | ug/L  | 10 U |
| bis(2-Chloroisopropyl) ether | 10    | ug/L  | 10 U |
| 4-Methylphenol               | 10    | ug/L  | 10 U |
| N-Nitrosodi-n-propylamine    | 10    | ug/L  | 10 U |
| Hexachloroethane             | 10    | ug/L  | 10 U |
| Nitrobenzene                 | 10    | ug/L  | 10 U |
| Isophorone                   | 10    | ug/L  | 10 U |
| 2-Nitrophenol                | 10    | ug/L  | 10 U |
| 2,4-Dimethylphenol           | 10    | ug/L  | 10 U |
| Benzoic acid                 | 50    | ug/L  | 50 U |
| bis(2-Chloroethoxy)methane   | 10    | ug/L  | 10 U |
| 2,4-Dichlorophenol           | 10    | ug/L  | 10 U |
| 1,2,4-Trichlorobenzene       | 10    | ug/L  | 10 U |
| Naphthalene                  | 10    | ug/L  | 10 U |
| 4-Chloroaniline              | 20    | ug/L  | 20 U |
| Hexachlorobutadiene          | 10    | ug/L  | 10 U |
| 4-Chloro-3-methylphenol      | 20    | ug/L  | 20 U |
| 2-Methylnaphthalene          | 10    | ug/L  | 10 U |
| Hexachlorocyclopentadiene    | 10    | ug/L  | 10 U |
| 2,4,6-Trichlorophenol        | 10    | ug/L  | 10 U |
| 2,4,5-Trichlorophenol        | 10    | ug/L  | 10 U |
| 2-Chloronaphthalene          | 10    | ug/L  | 10 U |
| 2-Nitroaniline               | 50    | ug/L  | 50 U |
| Dimethyl phthalate           | 10    | ug/L  | 10 U |
| Acenaphthylene               | 10    | ug/L  | 10 U |
| 2,6-Dinitrotoluene           | 10    | ug/L  | 10 U |
| 3-Nitroaniline               | 50    | ug/L  | 50 U |
| Acenaphthene                 | 10    | ug/L  | 10 U |
| 2,4-Dinitrophenol            | 50    | ug/L  | 50 U |
| 4-Nitrophenol                | 50    | ug/L  | 50 U |
| Dibenzofuran                 | 10    | ug/L  | 10 U |
| 2,4-Dinitrotoluene           | 10    | ug/L  | 10 U |
| Diethyl phthalate            | 10    | ug/L  | 10 U |
| 4-Chlorophenyl phenyl ether  | 10    | ug/L  | 10 U |
| Fluorene                     | 10    | ug/L  | 10 U |
| 4-Nitroaniline               | 50    | ug/L  | 50 U |
| 4,6-Dinitro-2-methylphenol   | 50    | ug/L  | 50 U |
| N-Nitrosodiphenylamine       | 10    | ug/L  | 10 U |
| 4-Bromophenyl phenyl ether   | 10    | ug/L  | 10 U |
| Hexachlorobenzene            | 10    | ug/L  | 10 U |

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 18, 1996

|                       |                       |             |
|-----------------------|-----------------------|-------------|
| EPA 8270B             | GTEL Sample Number    | M6110437-01 |
| Semivolatile Organics | Client ID             | FIELD BLANK |
| Matrix: Aqueous       | Date Sampled          | 11/20/96    |
|                       | Date Prepared         | 11/26/96    |
|                       | Date Analyzed         | 12/03/96    |
|                       | Adjustment Multiplier | 1.00        |

## Reporting

| Analyte                     | Limit | Units |      |
|-----------------------------|-------|-------|------|
| Pentachlorophenol           | 50    | ug/L  | 50 U |
| Phenanthrene                | 10    | ug/L  | 10 U |
| Anthracene                  | 10    | ug/L  | 10 U |
| Di-n-butyl phthalate        | 10    | ug/L  | 10 U |
| Fluoranthene                | 10    | ug/L  | 10 U |
| Pyrene                      | 10    | ug/L  | 10 U |
| Butylbenzyl phthalate       | 10    | ug/L  | 10 U |
| 3,3'-Dichlorobenzidine      | 20    | ug/L  | 20 U |
| Benzo[a]anthracene          | 10    | ug/L  | 10 U |
| Chrysene                    | 10    | ug/L  | 10 U |
| bis(2-Ethylhexyl) phthalate | 10    | ug/L  | 10 U |
| Di-n-octyl phthalate        | 10    | ug/L  | 10 U |
| Benzo[b]fluoranthene        | 10    | ug/L  | 10 U |
| Benzo[k]fluoranthene        | 10    | ug/L  | 10 U |
| Benzo[a]pyrene              | 10    | ug/L  | 10 U |
| Indeno[1,2,3-cd]pyrene      | 10    | ug/L  | 10 U |
| Dibenzo[a,h]anthracene      | 10    | ug/L  | 10 U |
| Benzo[g,h,i]perylene        | 10    | ug/L  | 10 U |
| Carbazole                   | 10    | ug/L  | 10 U |

Login Number: M6110437

Project ID (number): 7094104115110

Project ID (name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 18, 1996

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Footnotes and Comments

Symbol keys (may appear beside values)

- \* - Indicates the analyte has been qualified in the narrative section of the report.
- d - Indicates the analyte was reported from a dilution other than that indicated on the report page.
- B - Organic Analyses - Indicates the analyte is found in the associated blank as well as in the sample.
- B - Inorganic Analyses - Indicates an estimated value below the EPA Contract Required Detection Limit.
- G - Indicates an estimated surrogate recovery due to dilutions.
- J - Indicates an estimated value below the reporting limit.
- U - Indicates the analyte was analyzed for but not detected.
- NA - Matrix Spike Results - Not Applicable, since the Sample Conc. exceeded four times the Spike Added.
- NA - Matrix Spike Duplicate RPD Results - Not Applicable, since the Sample Conc. exceeded four times the Spike Added.
- NA - Serial Dilution RPD Results - Not Applicable, since the Sample Conc. was less than  
five times the CLP Contract Required Detection Limit.

Semi-Volatile Organics

Method: EPA 8270B

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", SW-846, Third Edition including promulgated Update II. Analyte list may be modified to include additional compounds.



GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 18, 1996

| EPA 8270B             | GTEL Sample Number    | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|-----------------------|-----------------------|-------------|------------------|-------------|-------------|
| Semivolatile Organics | Client ID             | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Matrix: Solids        | Date Sampled          | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
|                       | Date Prepared         | 12/03/96    | 12/03/96         | 12/03/96    | 12/03/96    |
|                       | Date Analyzed         | 12/05/96    | 12/05/96         | 12/05/96    | 12/05/96    |
|                       | Adjustment Multiplier | 1.00        | 1.00             | 1.00        | 1.00        |
|                       | Percent Solids        | 55.0        | 51.9             | 62.4        | 72.9        |

## Reporting

| Analyte                      | Limit | Units | Concentration: Dry Weight |        |        |        |
|------------------------------|-------|-------|---------------------------|--------|--------|--------|
| Phenol                       | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| bis(2-Chloroethyl) ether     | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2-Chlorophenol               | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 1,3-Dichlorobenzene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 1,4-Dichlorobenzene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Benzyl alcohol               | 660   | ug/kg | 660 U                     | 660 U  | 660 U  | 660 U  |
| 1,2-Dichlorobenzene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2-Methylphenol               | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| bis(2-Chloroisopropyl) ether | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 4-Methylphenol               | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| N-Nitrosodi-n-propylamine    | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Hexachloroethane             | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Nitrobenzene                 | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Isophorone                   | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2-Nitrophenol                | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2,4-Dimethylphenol           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Benzoic acid                 | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| bis(2-Chloroethoxy)methane   | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2,4-Dichlorophenol           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 1,2,4-Trichlorobenzene       | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Naphthalene                  | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 4-Chloroaniline              | 660   | ug/kg | 660 U                     | 660 U  | 660 U  | 660 U  |
| Hexachlorobutadiene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 4-Chloro-3-methylphenol      | 660   | ug/kg | 660 U                     | 660 U  | 660 U  | 660 U  |
| 2-Methylnaphthalene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Hexachlorocyclopentadiene    | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2,4,6-Trichlorophenol        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2,4,5-Trichlorophenol        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2-Chloronaphthalene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2-Nitroaniline               | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| Dimethyl phthalate           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Acenaphthylene               | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 2,6-Dinitrotoluene           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 3-Nitroaniline               | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| Acenaphthene                 | 330   | ug/kg | 330 U                     | 330 U  | 160 J  | 330 U  |
| 2,4-Dinitrophenol            | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| 4-Nitrophenol                | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| Dibenzofuran                 | 330   | ug/kg | 330 U                     | 330 U  | 84 J   | 330 U  |
| 2,4-Dinitrotoluene           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Diethyl phthalate            | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 4-Chlorophenyl phenyl ether  | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Fluorene                     | 330   | ug/kg | 330 U                     | 330 U  | 130 J  | 330 U  |
| 4-Nitroaniline               | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| 4,6-Dinitro-2-methylphenol   | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| N-Nitrosodiphenylamine       | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 4-Bromophenyl phenyl ether   | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 18, 1996

| EPA 8270B             | GTEL Sample Number    | M6110437-07 | M6110437-08      | M6110437-09 | M6110437-10 |
|-----------------------|-----------------------|-------------|------------------|-------------|-------------|
| Semivolatile Organics | Client ID             | RP1-4       | POND 1 DUPLICATE | RP1-5       | RP2-4       |
| Matrix: Solids        | Date Sampled          | 11/20/96    | 11/20/96         | 11/21/96    | 11/21/96    |
|                       | Date Prepared         | 12/03/96    | 12/03/96         | 12/03/96    | 12/03/96    |
|                       | Date Analyzed         | 12/05/96    | 12/05/96         | 12/05/96    | 12/05/96    |
|                       | Adjustment Multiplier | 1.00        | 1.00             | 1.00        | 1.00        |
|                       | Percent Solids        | 55.0        | 51.9             | 62.4        | 72.9        |

## Reporting

| Analyte                     | Limit | Units | Concentration: Dry Weight |        |        |        |
|-----------------------------|-------|-------|---------------------------|--------|--------|--------|
| Hexachlorobenzene           | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Pentachlorophenol           | 1700  | ug/kg | 1700 U                    | 1700 U | 1700 U | 1700 U |
| Phenanthrene                | 330   | ug/kg | 330 U                     | 330 U  | 300 J  | 87 J   |
| Anthracene                  | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Di-n-butyl phthalate        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Fluoranthene                | 330   | ug/kg | 330 U                     | 330 U  | 170 J  | 160 J  |
| Pyrene                      | 330   | ug/kg | 330 U                     | 330 U  | 130 J  | 150 J  |
| Butylbenzyl phthalate       | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| 3,3'-Dichlorobenzidine      | 660   | ug/kg | 660 U                     | 660 U  | 660 U  | 660 U  |
| Benzo[a]anthracene          | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 80 J   |
| Chrysene                    | 330   | ug/kg | 330 U                     | 330 U  | 55 J   | 100 J  |
| bis(2-Ethylhexyl) phthalate | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Di-n-octyl phthalate        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Benzo[b]fluoranthene        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 88 J   |
| Benzo[k]fluoranthene        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 57 J   |
| Benzo[a]pyrene              | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 81 J   |
| Indeno[1,2,3-cd]pyrene      | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 47 J   |
| Dibenzo[a,h]anthracene      | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |
| Benzo[g,h,i]perylene        | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 60 J   |
| Carbazole                   | 330   | ug/kg | 330 U                     | 330 U  | 330 U  | 330 U  |

|                       |                       |             |
|-----------------------|-----------------------|-------------|
| EPA 8270B             | GTEL Sample Number    | M6110437-11 |
| Semivolatile Organics | Client ID             | RP2-5       |
| Matrix: Solids        | Date Sampled          | 11/21/96    |
|                       | Date Prepared         | 12/03/96    |
|                       | Date Analyzed         | 12/05/96    |
|                       | Adjustment Multiplier | 1.00        |
|                       | Percent Solids        | 80.2        |

## Reporting

| Analyte                      | Limit | Units | Concentration: Dry Weight |
|------------------------------|-------|-------|---------------------------|
| Phenol                       | 330   | ug/kg | 330 U                     |
| bis(2-Chloroethyl) ether     | 330   | ug/kg | 330 U                     |
| 2-Chlorophenol               | 330   | ug/kg | 330 U                     |
| 1,3-Dichlorobenzene          | 330   | ug/kg | 330 U                     |
| 1,4-Dichlorobenzene          | 330   | ug/kg | 330 U                     |
| Benzyl alcohol               | 660   | ug/kg | 660 U                     |
| 1,2-Dichlorobenzene          | 330   | ug/kg | 330 U                     |
| 2-Methylphenol               | 330   | ug/kg | 330 U                     |
| bis(2-Chloroisopropyl) ether | 330   | ug/kg | 330 U                     |
| 4-Methylphenol               | 330   | ug/kg | 330 U                     |
| N-Nitrosodi-n-propylamine    | 330   | ug/kg | 330 U                     |
| Hexachloroethane             | 330   | ug/kg | 330 U                     |
| Nitrobenzene                 | 330   | ug/kg | 330 U                     |
| Isophorone                   | 330   | ug/kg | 330 U                     |

GTEL Client ID: 7094104115110  
 Login Number: M6110437  
 Project ID (Number): 7094104115110  
 Project ID (Name): RFI UNIVERSITY OF MARYLAND

## ANALYTICAL RESULTS

Date of Report: Dec 18, 1996

EPA 8270B GTEL Sample Number M6110437-11  
 Semivolatile Organics Client ID RP2-5  
 Matrix: Solids Date Sampled 11/21/96  
 Date Prepared 12/03/96  
 Date Analyzed 12/05/96  
 Adjustment Multiplier 1.00  
 Percent Solids 80.2

## Reporting

| Analyte                     | Limit | Units | Concentration: Dry Weight |
|-----------------------------|-------|-------|---------------------------|
| 2-Nitrophenol               | 330   | ug/kg | 330 U                     |
| 2,4-Dimethylphenol          | 330   | ug/kg | 330 U                     |
| Benzoic acid                | 1700  | ug/kg | 1700 U                    |
| bis(2-Chloroethoxy)methane  | 330   | ug/kg | 330 U                     |
| 2,4-Dichlorophenol          | 330   | ug/kg | 330 U                     |
| 1,2,4-Trichlorobenzene      | 330   | ug/kg | 330 U                     |
| Naphthalene                 | 330   | ug/kg | 330 U                     |
| 4-Chloroaniline             | 660   | ug/kg | 660 U                     |
| Hexachlorobutadiene         | 330   | ug/kg | 330 U                     |
| 4-Chloro-3-methylphenol     | 660   | ug/kg | 660 U                     |
| 2-Methylnaphthalene         | 330   | ug/kg | 330 U                     |
| Hexachlorocyclopentadiene   | 330   | ug/kg | 330 U                     |
| 2,4,6-Trichlorophenol       | 330   | ug/kg | 330 U                     |
| 2,4,5-Trichlorophenol       | 330   | ug/kg | 330 U                     |
| 2-Chloronaphthalene         | 330   | ug/kg | 330 U                     |
| 2-Nitroaniline              | 1700  | ug/kg | 1700 U                    |
| Dimethyl phthalate          | 330   | ug/kg | 330 U                     |
| Acenaphthylene              | 330   | ug/kg | 330 U                     |
| 2,6-Dinitrotoluene          | 330   | ug/kg | 330 U                     |
| 3-Nitroaniline              | 1700  | ug/kg | 1700 U                    |
| Acenaphthene                | 330   | ug/kg | 330 U                     |
| 2,4-Dinitrophenol           | 1700  | ug/kg | 1700 U                    |
| 4-Nitrophenol               | 1700  | ug/kg | 1700 U                    |
| Dibenzofuran                | 330   | ug/kg | 330 U                     |
| 2,4-Dinitrotoluene          | 330   | ug/kg | 330 U                     |
| Diethyl phthalate           | 330   | ug/kg | 330 U                     |
| 4-Chlorophenyl phenyl ether | 330   | ug/kg | 330 U                     |
| Fluorene                    | 330   | ug/kg | 330 U                     |
| 4-Nitroaniline              | 1700  | ug/kg | 1700 U                    |
| 4,6-Dinitro-2-methylphenol  | 1700  | ug/kg | 1700 U                    |
| N-Nitrosodiphenylamine      | 330   | ug/kg | 330 U                     |
| 4-Bromophenyl phenyl ether  | 330   | ug/kg | 330 U                     |
| Hexachlorobenzene           | 330   | ug/kg | 330 U                     |
| Pentachlorophenol           | 1700  | ug/kg | 1700 U                    |
| Phenanthrene                | 330   | ug/kg | 330 U                     |
| Anthracene                  | 330   | ug/kg | 330 U                     |
| Di-n-butyl phthalate        | 330   | ug/kg | 330 U                     |
| Fluoranthene                | 330   | ug/kg | 330 U                     |
| Pyrene                      | 330   | ug/kg | 330 U                     |
| Butylbenzyl phthalate       | 330   | ug/kg | 330 U                     |
| 3,3'-Dichlorobenzidine      | 660   | ug/kg | 660 U                     |
| Benzo[a]anthracene          | 330   | ug/kg | 330 U                     |
| Chrysene                    | 330   | ug/kg | 330 U                     |
| bis(2-Ethylhexyl) phthalate | 330   | ug/kg | 330 U                     |
| Di-n-octyl phthalate        | 330   | ug/kg | 330 U                     |
| Benzo[b]fluoranthene        | 330   | ug/kg | 330 U                     |

GTEL Client ID: 7094104115110

## ANALYTICAL RESULTS

Login Number: M6110437

Project ID (Number): 7094104115110

Project ID (Name): RFI UNIVERSITY OF MARYLAND

Date of Report: Dec 18, 1996

|                       |                       |             |
|-----------------------|-----------------------|-------------|
| EPA 8270B             | GTEL Sample Number    | M6110437-11 |
| Semivolatile Organics | Client ID             | RP2-5       |
| Matrix: Solids        | Date Sampled          | 11/21/96    |
|                       | Date Prepared         | 12/03/96    |
|                       | Date Analyzed         | 12/05/96    |
|                       | Adjustment Multiplier | 1.00        |
|                       | Percent Solids        | 80.2        |

| Analyte                | Limit | Units | Concentration: Dry Weight |
|------------------------|-------|-------|---------------------------|
| Benzo[k]fluoranthene   | 330   | ug/kg | 330 U                     |
| Benzo[a]pyrene         | 330   | ug/kg | 330 U                     |
| Indeno[1,2,3-cd]pyrene | 330   | ug/kg | 330 U                     |
| Dibenzo[a,h]anthracene | 330   | ug/kg | 330 U                     |
| Benzo[g,h,i]perylene   | 330   | ug/kg | 330 U                     |
| Carbazole              | 330   | ug/kg | 330 U                     |

## **APPENDIX D**

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### **NEI/GTEL CORRESPONDENCE**



TOTAL ANALYTICAL SERVICES FOR A SAFE ENVIRONMENT

nytest environmental inc.

Buchart Horn, Inc.  
The Industrial Plaza of York  
445 West Philadelphia Street  
P.O. Box 15040  
York, PA 17405-7040

Attention: Randy Deardorf

Subject: Herbicide Analysis for the University of Maryland Project,  
Phase I, RFI

Dear Mr. Deardorf:

This letter is provided as a follow-up in regard to our conversation concerning the fact that Nytest Environmental, Inc. did not analyze for diallate, requested as a herbicide, by method 8150, for the referenced project.

On November 11, 1996, I received paperwork from you detailing the final analytical requirements and sample numbers for the referenced project. Diallate was listed with the herbicide list and I mistakenly interpreted this compound to be another name for 2,4-D and therefore did not request the laboratory to run for it. Once you received the data, it was discovered this compound was not reported and it is out of holding time at this time.

Once the error was identified, I further researched this compound and I hope your project was not significantly impacted in a negative way. Here is the information, I have found concerning diallate:

- . It is a compound that would require a separate run and some research to accommodate.
- . It is not a herbicide that can be run by 8150.
- . Our chemical directory lists diallate as a pesticide, carbamate, as follows:  
S-(2-3-dichloroallyl)N-N'-diisopropyl thiocarbamate, Avadex, CAS#: 2303-16-4.  
This compound is a different acid group than compounds associated with method 8150.
- . I contacted Core Laboratories in Indiana, one of the few labs who analyze for carbamates in drinking water by method 531.1 but they do not run diallate by this method or any other method.

box 1518 ☐ 60 seaview blvd., port washington, ny 11050 ☐ (516) 625-5500  
fax (516) 625-1274

## nytest environmental inc.

- . I contacted Lancaster Laboratories who analyze s for several more of the drinking water methods for the Synthetic Organic Compounds which have more extensive pesticide, herbicide and semivolatile lists. These methods are not always applicable for soils and Lancaster has not run for diallate nor is it detected or listed as a compound in the following methods they do run: 505 (includes some pesticides, PCBs and semivolatiles), method 515.1 (includes some herbicides and pesticides), 525.1 (includes some pesticides and semivolatiles) and 531.1 (carbamates).
- . Diallate is a carbamate which are usually run utilizing HPLC. Both of the referenced laboratories have HPLC capabilities but do not run for diallate.
- . Diallate is listed as a semivolatile (method 8270) on the Appendix IX list which is the list of parameters normally tested to determine leachate from landfills and is not always a routinely run list. Diallate is not a TCL or PP semivolatile and is not run at NEI/GTEL in New Hampshire. To determine if this laboratory could run this compound by 8270, I should have given the laboratory advance notice. This would have taken some research and development, and could have potentially not worked..
- . Diallate is identified as a U listed waste, U-062, but I have been unable to identify any concentration constituent in waste which may mean there is no standard for this compound.
- . One of our chemist found a reference that diallate may potentially be able to be run by method 8080/8081 which is routinely a pesticide/PCB method. Again this is a nonstandard compound for this method, would require research and development and may potentially not work.

I hope this letter provides some useful information. I regret that I overlooked this compound up front when perhaps NEI could have done the research and development or we may have determined along with the EPA that this compound could be dropped from the requirements. Research and development and an additional method could have had a cost impact on the project. NEI in New York has run the Semivolatile Appendix IX list by Method 8270 which may be the approach to take if this compound must be run but we would need to order the standard.

NEI and I apologize for any inconvenience we may have caused you and your client. Please do not hesitate to call me at (609) 829-7390 if we may need to accommodate this project requirement or lend further assistance .

Sincerely,

Nytest Environmental Inc.

  
Anne W. Lee  
Account Executive